10<sup>756214</sup>9 JC10 Rec'd PCT/PTO 23 DEC 2005

### Appendix A

#### Claim Amendments

1. (Currently amended) Compounds A compound of formula I

**(l)** 

in which

- R1 is halogen, nitro, amino, mono- or di-1-4C-alkylamino, 1-4C-alkyl, hydroxyl, 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, 3-7C-cycloalkoxy, 3-7C-cycloalkylmethoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy,
- R2 is hydrogen, halogen or 1-4C-alkoxy, and
- R3 is hydrogen or 1-4C-alkoxy, or
- R2 and R3 bound to the benzo ring moiety in ortho-position to each other together form a 1-2C-alkylenedioxy bridge,

- R2 and R3 bound to the benzo ring moiety in ortho-position to each other together form a completely or predominantly fluorine-substituted 1-2C-alkylenedioxy bridge, or
- R1 and R2 bound to the benzo ring moiety in ortho-position to each other together form a 1-2C-alkylenedioxy bridge and R3 is hydrogen, or
- R1 and R2 bound to the benzo ring moiety in ortho-position to each other together form a completely or predominantly fluorine-substituted 1-2C-alkylenedioxy bridge and R3 is hydrogen,
- R4 is hydrogen, fluorine, chlorine, 1-4C-alkyl, trifluoromethyl, cyclopropyl, cyano, 1-4C-alkoxycarbonyl or -CH2-O-R411, in which
- R411 is hydrogen, 1-4C-alkyl, 1-4C-alkoxy-2-4-alkyl or 1-4C-alkylcarbonyl,

R41 is hydrogen or 1-4C-alkyl,

R5 is hydrogen, and

R51 is hydrogen,

or

R4 is hydrogen, fluorine, chlorine or 1-4C-alkyl,

R41 is hydrogen or 1-4C-alkyl,

R5 is hydrogen, and

- R51 is hydrogen,
- R6 is 1-6C-alkyl, amino, formyl, or 1-4C-alkyl substituted by R61, in which
- R61 is 1-4C-alkoxycarbonyl, carboxyl, 1-4C-alkoxy, hydroxyl, halogen or -N(R611)R612, in which
- R611 is hydrogen, 1-4C-alkyl, 3-7C-cycloalkyl or 3-7C-cycloalkyl-1-4C-alkyl, and
- R612 is hydrogen or 1-4C-alkyl, or
- R611 and R612 together and with inclusion of the nitrogen atom to which they are bound form a radical Het1, in which
- Hetl is a 5- to 7-membered saturated heterocyclic ring radical comprising one nitrogen atom, to which R611 and R612 are bound, and, optionally, one further heteroatom selected from [[a]] the group consisting of nitrogen, oxygen and sulfur, and optionally substituted by R613 on a ring nitrogen atom, in which
- R613 is 1-4C-alkyl, 3-7C-cycloalkyl, 3-7C-cycloalkyl-1-4C-alkyl, hydroxy-2-4C-alkyl, 1-4C-alkoxy-2-4C-alkyl, amino-2-4C-alkyl, mono- or di-1-4C-alkylamino-2-4C-alkyl, formyl, pyridyl or pyrimidinyl,

- R7 is phenyl, Het2, R71- and/or R72- and/or R73-substituted phenyl, R74- and/or R75-substituted Het2, naphthyl, or R76- and/or R77-substituted naphthyl, in which
- Het2 is a monocyclic or fused bicyclic 5 to 10-membered heteroaryl radical comprising one to three heteroatoms, each of which is selected from [[a]] the group consisting of nitrogen, oxygen and sulfur,
- R71 is hydroxyl, halogen, nitro, cyano, trifluoromethyl, 1-4C-alkyl, 1-4C-alkoxy, amino, mono- or di-1-4C-alkylamino, 1-4C-alkylsulphonylamino, arylsulphonylamino, 1-4C-alkoxycarbonyl, carboxyl, 1-4C-alkylthio, aryloxy-2-4C-alkoxy, aryloxy-1-4C-alkyl, aryloxy, aryl-1-4C-alkoxy, aryl, 1-4C-alkoxy-2-4C-alkoxy, 1-4C-alkoxy-1-4C-alkyl, hydroxy-2-4C-alkoxy, amino-2-4C-alkoxy, mono- or di-1-4C-alkylamino-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy, in which
- aryl is phenyl or R711-substituted phenyl, in which
  R711 is halogen, 1-4C-alkyl, 1-4C-alkoxy, nitro or cyano,
  R72 is halogen, 1-4C-alkyl, 1-4C-alkoxy or 1-4Calkoxycarbonyl,
- R73 is 1-4C-alkyl or 1-4C-alkoxy,

- R74 is halogen, 1-4C-alkyl, trifluoromethyl, 1-4C-alkoxy, cyano, amino, mono- or di-1-4C-alkylamino, 1-4C-alkoxycarbonyl, morpholino, carboxyl, nitro, phenyl or phenyloxy,
- R75 is 1-4C-alkyl or halogen,
- R76 is halogen, hydroxyl, 1-4C-alkyl, 1-4C-alkoxy, carboxyl or 1-4C-alkoxycarbonyl,
- R77 is 1-4C-alkyl or 1-4C-alkoxy,
- R8 is phenyl, phenylcarbonyl, -C(O)-N(R82)R83 or -C(O)-OR9, in which
- R82 is hydrogen, 1-4C-alkyl, 3-7C-cycloalkyl, 3-7C-cycloalkyl-1-4C-alkyl, phenyl or phenyl-1-4C-alkyl, and R83 is hydrogen or 1-4C-alkyl, or
- R82 and R83 together and with inclusion of the nitrogen atom, to which they are bound, form a heterocyclic ring radical selected from the group consisting of pyrrolidinyl, piperidinyl, morpholinyl [[or]] and N-(1-4C-alkyl)-piperazinyl,
- R9 is 1-4C-alkyl, [[;]]
- or a salt, stereoisomer, hydrate or hydrate of a salt thereof;

under the provisio, proviso that this subgroup of compounds of formula I, wherein the combination of all of the

following restrictions a.) to c.) apply, is thereof disclaimed:

a.) the substitution pattern of the left Ri- and/or R2and/or R3-substituted benzo ring of the dihydroisoquinoline moiety of the pyrrolodihydroisoquinoline scaffold shown in formula I is as follows:

in which

R' and R'' can be bonded at any possible position of the benzo ring, and

R' is hydroxyl, 1-4C-alkoxy or trifluoromethoxy,
R''is hydrogen or 1-4C-alkoxy,

or R' and R'' bound to the benzo ring moiety in orthoposition to each other together form a 1-2C-alkylenedioxy bridge,

and

b.) R4 is hydrogen, andR41 is hydrogen, andR5 is hydrogen, andR51 is hydrogen,

and

- c.) R8 is -C(0)-OR9, in which
   R9 is 1-4C-alkyl+
- and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.
- 2. (Currently amended) Compounds A compound of formula I according to claim 1, in which
- R1 is halogen, nitro, amino, 1-4C-alkyl, 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy,
- R2 is 1-4C-alkoxy,
- R3 is hydrogen,
- and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1-a] isoquinoline ring,
- R4 is hydrogen or 1-4C-alkyl,
- R41 is hydrogen or 1-4C-alkyl,
- R5 is hydrogen,
- R51 is hydrogen,
- R6 is 1-6C-alkyl, formyl, or 1-4C-alkyl substituted by R61, in which
- R61 is 1-4C-alkoxycarbonyl,

- R7 is phenyl, Het2, R71- and/or R72- and/or R73-substituted phenyl, R74- and/or R75-substituted Het2, or naphthyl, in which
- Het2 is a heteroaryl radical selected from the group consisting of furanyl, thiophenyl, pyrrolyl, pyridinyl, quinolyl, indolyl, benzothiophenyl and benzofuranyl,
- R71 is hydroxyl, 1-4C-alkoxy, amino or mono- or di-1-4C-alkylamino,
- R72 is 1-4C-alkyl or 1-4C-alkoxy,
- R73 is 1-4C-alkyl or 1-4C-alkoxy,
- R74 is 1-4C-alkyl, trifluoromethyl, 1-4C-alkoxy, 1-4C-alkoxycarbonyl, nitro, phenyl or phenyloxy,
- R75 is 1-4C-alkyl,
- R8 is phenyl, phenylcarbonyl, or -C(O)-N(R82)R83, in which
- R82 is hydrogen, 1-4C-alkyl, 3-7C-cycloalkyl or phenyl, and
- R83 is hydrogen or 1-4C-alkyl, or
- R82 and R83 together and with inclusion of the nitrogen atom, to which they are bound, form a heterocyclic ring radical selected from the group consisting of pyrrolidinyl and piperidinyl,
- or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

- 3. (Currently amended) Compounds A compound of formula I according to claim 1, in which
- R1 is halogen, nitro, amino, 1-4C-alkyl, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy,

with the provisio that R1 is not trifluoromethoxy,

- R2 is 1-4C-alkoxy, and
- R3 is hydrogen, or
- R1 and R2 bound to the benzo ring moiety in ortho-position to each other together form a completely or predominantly fluorine-substituted 1-2C-alkylenedioxy bridge and R3 is hydrogen,
- and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1-a] isoquinoline ring,
- R4 is hydrogen or 1-4C-alkyl,
- R41 is hydrogen or 1-4C-alkyl,
- R5 is hydrogen,
- R51 is hydrogen,

- R6 is 1-6C-alkyl, formyl, or 1-4C-alkyl substituted by R61, in which
- R61 is 1-4C-alkoxycarbonyl,
- R7 is phenyl, Het2, R71- and/or R72- and/or R73-substituted phenyl, R74- and/or R75-substituted Het2, or naphthyl, in which
- Het2 is a heteroaryl radical selected from the group consisting of furanyl, thiophenyl, pyrrolyl, pyridinyl, quinolyl, indolyl, benzothiophenyl and benzofuranyl,
- R71 is hydroxyl, 1-4C-alkoxy, amino or mono- or di-1-4C-alkylamino,
- R72 is 1-4C-alkyl or 1-4C-alkoxy,
- R73 is 1-4C-alkyl or 1-4C-alkoxy,
- R74 is 1-4C-alkyl, trifluoromethyl, 1-4C-alkoxy, 1-4C-alkoxy, nitro, phenyl or phenyloxy,
- R75 is 1-4C-alkyl,
- R8 is phenyl, phenylcarbonyl, -C(O)-N(R82)R83 or -C(O)-OR9, in which
- R82 is hydrogen, 1-4C-alkyl, 3-7C-cycloalkyl or phenyl, and
- R83 is hydrogen or 1-4C-alkyl, or
- R82 and R83 together and with inclusion of the nitrogen atom, to which they are bound, form a heterocyclic ring

radical selected from the group consisting of pyrrolidinyl and piperidinyl,

R9 is 1-4C-alkyl,

# or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

- 4. (Currently amended) Compounds A compound of formula I according to claim 1, in which
- R1 is halogen, nitro, amino, 1-4C-alkyl, 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy,
- R2 is 1-4C-alkoxy,
- R3 is 1-4C-alkoxy,
- and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1 a]isoquinoline
  pyrrolo[2,1-a]isoquinoline
  ring,
- R4 is hydrogen or 1-4C-alkyl,
- R41 is hydrogen or 1-4C-alkyl,
- R5 is hydrogen,
- R51 is hydrogen,

- R6 is 1-6C-alkyl, formyl, or 1-4C-alkyl substituted by R61, in which
- R61 is 1-4C-alkoxycarbonyl,
- R7 is phenyl, Het2, R71- and/or R72- and/or R73-substituted phenyl, R74- and/or R75-substituted Het2, or naphthyl, in which
- Het2 is a heteroaryl radical selected from the group consisting of furanyl, thiophenyl, pyrrolyl, pyridinyl, quinolyl, indolyl, benzothiophenyl and benzofuranyl,
- R71 is hydroxyl, 1-4C-alkoxy, amino or mono- or di-1-4C-alkylamino,
- R72 is 1-4C-alkyl or 1-4C-alkoxy,
- R73 is 1-4C-alkyl or 1-4C-alkoxy,
- R74 is 1-4C-alkyl, trifluoromethyl, 1-4C-alkoxy, 1-4C-alkoxycarbonyl, nitro, phenyl or phenyloxy,
- R75 is 1-4C-alkyl,
- R8 is phenyl, phenylcarbonyl, -C(O)-N(R82)R83 or -C(O)-OR9, in which
- R82 is hydrogen, 1-4C-alkyl, 3-7C-cycloalkyl or phenyl, and
- R83 is hydrogen or 1-4C-alkyl, or
- R82 and R83 together and with inclusion of the nitrogen atom, to which they are bound, form a heterocyclic ring

radical selected from the group consisting of pyrrolidinyl and piperidinyl,

R9 is 1-4C-alkyl,

# or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

- 5. (Currently amended) Compounds A compound of formula I according to claim 1, in which
- R1 is halogen, nitro, amino, 1-4C-alkyl, 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy,
- R2 is 1-4C-alkoxy,
- R3 is hydrogen,
- and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1 a]isoquinoline
  pyrrolo[2,1-a]isoquinoline ring,

R4 is 1-4C-alkyl,

R41 is hydrogen or 1-4C-alkyl,

R5 is hydrogen,

R51 is hydrogen,

- R6 is 1-6C-alkyl, formyl, or 1-4C-alkyl substituted by R61, in which
- R61 is 1-4C-alkoxycarbonyl,
- R7 is phenyl, Het2, R71- and/or R72- and/or R73-substituted phenyl, R74- and/or R75-substituted Het2, or naphthyl, in which
- Het2 is a heteroaryl radical selected from the group consisting of furanyl, thiophenyl, pyrrolyl, pyridinyl, quinolyl, indolyl, benzothiophenyl and benzofuranyl,
- R71 is hydroxyl, 1-4C-alkoxy, amino or mono- or di-1-4C-alkylamino,
- R72 is 1-4C-alkyl or 1-4C-alkoxy,
- R73 is 1-4C-alkyl or 1-4C-alkoxy,
- R74 is 1-4C-alkyl, trifluoromethyl, 1-4C-alkoxy, 1-4C-alkoxy, nitro, phenyl or phenyloxy,
- R75 is 1-4C-alkyl,
- R8 is phenyl, phenylcarbonyl, -C(O)-N(R82)R83 or -C(O)-OR9, in which
- R82 is hydrogen, 1-4C-alkyl, 3-7C-cycloalkyl or phenyl, and
- R83 is hydrogen or 1-4C-alkyl, or
- R82 and R83 together and with inclusion of the nitrogen atom, to which they are bound, form a heterocyclic ring

radical selected from the group consisting of pyrrolidinyl and piperidinyl,

R9 is 1-4C-alkyl,

# or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

6. (Currently amended) Compounds A compound of formula I according to claim 1,

in which, in a first embodiment,

R1 is halogen, nitro, amino, mono- or di-1-4C-alkylamino, 1-4C-alkyl, 1-4C-alkoxy-2-4C-alkoxy, 3-7C-cycloalkoxy, 3-7C-cycloalkylmethoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy,

R2 is 1-4C-alkoxy,

R3 is hydrogen,

and none of R1 and R2 is bound to the 10-position of the pyrrolo[2.1-a]isoquinoline pyrrolo[2,1-a]-isoquinoline
ring,

and

R4 is hydrogen,

R41 is hydrogen,

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R5 is hydrogen, and
R51 is hydrogen;
or in which, in a second embodiment,
R1 is 1-4C-alkoxy,
   is 1-4C-alkoxy,
R2
R3
   is hydrogen,
and none of R1 and R2 is bound to the 10-position of the
   pyrrolo[2.1 a] isoquinoline
pyrrolo[2,1-a] isoquinoline
   ring,
and
R4 is 1-4C-alkyl,
R41 is hydrogen or 1-4C-alkyl,
R5 is hydrogen, and
R51 is hydrogen;
R6 is 1-6C-alkyl, or 1-4C-alkyl substituted by R61,
   which
R61 is 1-4C-alkoxycarbonyl;
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R7 is Het2, 4-hydroxy-3,5-dimethylphenyl, 3-dimethylamino-

phenyl, or naphthyl, in which

- Het2 is a monocyclic or fused bicyclic 5- to 10-membered heteroaryl radical comprising one to three heteroatoms, each of which is selected from [[a]] the group consisting of nitrogen, oxygen and sulfur,
- such as, for example, quinolyl, e.g. quinolin-4-yl;
- R8 is phenyl, phenylcarbonyl, -C(O)-N(R82)R83 or -C(O)-OR9, in which
- R82 is hydrogen, 1-4C-alkyl, 3-7C-cycloalkyl or phenyl, and R83 is hydrogen or 1-4C-alkyl, or
- R82 and R83 together and with inclusion of the nitrogen atom, to which they are bound, form a heterocyclic ring radical selected from the group consisting of pyrrolidinyl and piperidinyl,
- R9 is 1-4C-alkyl;
- or a salt, stereoisomer, hydrate or hydrate of a salt thereof
- and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.
- 7. (Currently amended)  $\frac{A}{A}$  compound of formula I according to claim 1,
- in which, in a first embodiment,

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is halogen, 1-4C-alkoxy-2-4C-alkoxy, or completely or
R1
   predominantly fluorine-substituted 1-4C-alkoxy,
   is 1-4C-alkoxy,
R3
   is hydrogen,
and none of R1 and R2 is bound to the 7- or 10-position of
    the
            pyrrolo[2.1-a]isoquinoline
                                            pyrrolo[2,1-a]-
    isoquinoline ring,
and
   is hydrogen,
R41 is hydrogen,
R5 is hydrogen, and
R51 is hydrogen;
or in which, in a second embodiment,
R1 is 1-4C-alkoxy,
   is 1-4C-alkoxy,
R2
   is hydrogen,
R3
and none of R1 and R2 is bound to the 7- or 10-position of
             pyrrolo[2.1 a]isoquinoline
                                             pyrrolo[2,1-a]-
    isoquinoline ring,
and
   is 1-4C-alkyl,
R41 is hydrogen or 1-4C-alkyl,
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R5 is hydrogen, and

R51 is hydrogen;

R6 is 1-4C-alkyl, or 1-4C-alkyl substituted by R61, in which

R61 is 1-4C-alkoxycarbonyl;

- R7 is Het2, 4-hydroxy-3,5-dimethylphenyl, 3-dimethylaminophenyl, or naphthyl, in which
- Het2 is a monocyclic or fused bicyclic 5- to 10-membered heteroaryl radical comprising one to three heteroatoms, each of which is selected from [[a]] the group consisting of nitrogen, oxygen and sulfur, such as, for example, quinolyl, e.g. quinolin 4 yl,

R8 is -C(O)-OR9, in which

R9 is 1-4C-alkyl;

- or a salt, stereoisomer, hydrate or hydrate of a salt thereof
- and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.
- 8. (Currently amended) Compounds A compound of formula I according to claim 1,

in which, in a first embodiment,
either

- R1 is halogen, 1-4C-alkyl, 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy,
- R2 is 1-4C-alkoxy, and
- R3 is 1-4C-alkoxy,

or

- R1 is 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy,
- R2 is halogen, and
- R3 is 1-4C-alkoxy,
- and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1-a]isoquinoline
  pyrrolo[2,1-a]isoquinoline ring;

or

either

- R1 is halogen, 1-4C-alkyl, 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy,
- R2 is 1-4C-alkoxy, and
- R3 is hydrogen,

or

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R1 is 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, or completely
    or predominantly fluorine-substituted 1-4C-alkoxy,
R2
   is halogen, and
R3 is hydrogen,
and none of R1 and R2 is bound to the 7- or 10-position of
            pyrrolo[2.1 a]isoquinoline
                                        pyrrolo[2,1-a]-
    isoquinoline ring;
and
R4 is 1-4C-alkyl,
R41 is hydrogen or 1-4C-alkyl,
R5 is hydrogen, and
R51 is hydrogen;
or in which, in a second embodiment,
R1 is 1-4C-alkoxy,
   is 1-4C-alkoxy,
R2
   is hydrogen,
R3
and none of R1 and R2 is bound to the 7- or 10-position of
   the
            pyrrolo[2.1-a]isoquinoline
                                             pyrrolo[2,1-a]-
   isoquinoline ring,
and
   is 1-4C-alkyl,
R4
R41 is hydrogen or 1-4C-alkyl,
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- R5 is hydrogen,
- R51 is hydrogen;
- R6 is 1-6C-alkyl, or 1-4C-alkyl substituted by R61, in which
- R61 is 1-4C-alkoxycarbonyl,
- R7 is phenyl, Het2, R71- and/or R72- and/or R73-substituted phenyl, R74- and/or R75-substituted Het2, or naphthyl, in which
- Het2 is a heteroaryl radical selected from the group consisting of furanyl, thiophenyl, pyrrolyl, pyridinyl, quinolyl, indolyl, benzothiophenyl and benzofuranyl,
- R71 is hydroxyl, 1-4C-alkoxy, amino or mono- or di-1-4C-alkylamino,
- R72 is 1-4C-alkyl or 1-4C-alkoxy,
- R73 is 1-4C-alkyl or 1-4C-alkoxy,
- R74 is 1-4C-alkyl, trifluoromethyl, 1-4C-alkoxy, 1-4C-alkoxy, nitro, phenyl or phenyloxy,
- R75 is 1-4C-alkyl,
- R8 is phenyl, phenylcarbonyl, -C(O)-N(R82)R83 or -C(O)-OR9, in which
- R82 is hydrogen, 1-4C-alkyl, 3-7C-cycloalkyl or phenyl, and
- R83 is hydrogen or 1-4C-alkyl, or

- R82 and R83 together and with inclusion of the nitrogen atom, to which they are bound, form a heterocyclic ring radical selected from the group consisting of pyrrolidinyl and piperidinyl,
- R9 is 1-4C-alkyl,
- or a salt, stereoisomer, hydrate or hydrate of a salt thereof
- and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.
- 9. (Currently amended) Compounds A compound of formula I according to claim 1,

in which

- R1 is halogen, 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy, such as e.g. chlorine, methoxy, 2 methoxy ethoxy or difluoromethoxy,
- R2 is 1-4C-alkoxy, such as-e.g. methoxy,
- R3 is 1-4C-alkoxy, such as e.g. methoxy,
- and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1-a]isoquinoline
  pyrrolo[2,1-a]isoquinoline ring,

and R4 is hydrogen, or 1-4C-alkyl such as e.g. methyl, R41 is hydrogen, or 1-4C-alkyl such as e.q. methyl, R5 is hydrogen, and R51 is hydrogen; or is halogen, 1-4C-alkoxy-2-4C-alkoxy, or completely or R1 predominantly fluorine-substituted 1-4C-alkoxy, such-as e.g. chlorine, 2 methoxy-ethoxy or difluoromethoxy, is 1-4C-alkoxy, such as e.g. methoxy, R2 R3 is hydrogen, and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1 a]isoquinoline pyrrolo[2,1-a]isoquinoline ring, and R4 is hydrogen, or 1-4C-alkyl such-as e.g. methyl, R41 is hydrogen, or 1-4C-alkyl such as e.g. methyl, R5 is hydrogen, and R51 is hydrogen; or

is 1-4C-alkoxy, such as e.g. methoxy,

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R2 is 1-4C-alkoxy, such as e.g. methoxy,
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R3 is hydrogen,

and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1-a]isoquinoline
pyrrolo[2,1-a]isoquinoline ring,

and

R4 is 1-4C-alkyl, such as e.g. methyl,

R41 is hydrogen, or 1-4C-alkyl such as e.g. methyl,

R5 is hydrogen, and

R51 is hydrogen;

R6 is 1-4C-alkyl, such as e.g. methyl;

R7 is 4-hydroxy-3,5-dimethylphenyl, 3-dimethylamino-phenyl, or

pyridyl, indolyl, thiophenyl, quinolinyl or naphthyl, such as e.g. pyridin 4 yl, indol 3 yl, thiophen 3 yl, quinolin 4 yl or naphthalen 1 yl,

R8 is -C(0) -OR9, in which

R9 is 1-2C-alkyl, such as e.g. ethyl,

or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

10. (Currently amended) Compounds A compound of formula I according to claim 1,

in which, in a first embodiment,

either

- R1 is bonded in the 8-position of the pyrrolo[2.1
  a]isoquinoline pyrrolo[2,1-a]-isoquinoline ring, and is
  chlorine, 2-methoxy-ethoxy or difluoromethoxy, and
- R2 is bonded in the 9-position of the pyrrolo[2.1a]isoquinoline pyrrolo[2,1-a]-isoquinoline ring, and is methoxy,

or

- R1 is bonded in the 9-position of the pyrrolo[2.1a]isoquinoline pyrrolo[2,1-a]-isoquinoline ring, and is
  chlorine, fluorine, methyl, nitro, amino or
  difluoromethoxy, and
- R2 is bonded in the 8-position of the pyrrolo[2.1a]isoquinoline pyrrolo[2,1-a]-isoquinoline ring, and is methoxy,

and

R3 is hydrogen,

and

either

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R4 is hydrogen,
R41 is hydrogen,
R5 is hydrogen, and
R51 is hydrogen,
or
R4
   is methyl,
R41 is hydrogen or methyl,
R5 is hydrogen, and
R51 is hydrogen;
or in which, in a second embodiment,
R1 is bonded in the 8-position of the pyrrolo[2.1
   a] isoquinoline pyrrolo[2,1-a] - isoquinoline ring, and is
   methoxy,
R2 is
        bonded in the 9-position of the pyrrolo[2.1-
   alisoquinoline pyrrolo[2,1-a]-isoquinoline ring, and is
   methoxy,
R3 is hydrogen,
and
R4 is methyl,
R41 is hydrogen or methyl,
R5 is hydrogen,
R51 is hydrogen;
```

- R6 is methyl or 2-methoxycarbonylethyl,
- R7 is 4-hydroxy-3,5-dimethylphenyl, 3-dimethylamino-phenyl, 3,4,5-trimethoxyphenyl, quinolinyl or naphthyl,
- R8 is phenylcarbonyl, -C(O)-N(R82)R83 or -C(O)-OR9, in which
- R82 is hydrogen, methyl, ethyl, iso-propyl, iso-butyl, cyclohexyl, cyclopropyl or phenyl, and
- R83 is hydrogen or methyl, or
- R82 and R83 together and with inclusion of the nitrogen atom, to which they are bound, form a pyrrolidinyl radical,
- R9 is methyl or ethyl,
- or a salt, stereoisomer, hydrate or hydrate of a salt thereof
- and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.
- 11. (Currently amended) Compounds A compound according to  $\frac{\text{claim 1}}{\text{claim 5}} \frac{\text{any of the claims 1 to 10}}{\text{to 10}}, \text{ wherein said compounds}$  have the formula I, in which

R1 is 1-4C-alkoxy, such as e.g. 1-2C alkoxy,

R2 is 1-4C-alkoxy, such as e.g. 1-2C-alkoxy,

R3 is hydrogen,

and none of R1 and R2 is bound to the 7- or 10-position of the pyrrolo[2.1-a] isoquinoline ring,

and in which

R4 is 1-4C-alkyl, such as e.g. methyl,

R41 is hydrogen, or 1-4C-alkyl such as e.g. methyl,

R5 is hydrogen,

R51 is hydrogen,

or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

12. (Currently amended) Compounds A compound according to claim 1 any of the claims 1 to 10, wherein said compounds have the formula I,

in which

either

R1 is halogen, nitro, amino, mono- or di-1-4C-alkylamino, 1-4C-alkyl, 1-4C-alkoxy-2-4C-alkoxy, 3-7C-cycloalkoxy,

```
3-7C-cycloalkylmethoxy, or completely or predominantly
    fluorine-substituted 1-4C-alkoxy, and
    is 1-4C-alkoxy,
R2
or
    is 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, 3-7C-
R1
    cycloalkoxy, 3-7C-cycloalkylmethoxy, or completely or
    predominantly fluorine-substituted 1-4C-alkoxy, and
R2 is halogen,
and
R3
    is hydrogen,
and none of R1 and R2 is bound to the 7- or 10-position of
    the
             pyrrolo[2.1-a] isoquinoline
                                              pyrrolo[2,1-a]-
    isoquinoline ring;
and in which
either
R4 is hydrogen,
R41 is hydrogen,
R5 is hydrogen, and
R51 is hydrogen,
or
R4
   is 1-4C-alkyl,
R41 is hydrogen or 1-4C-alkyl,
R5 is hydrogen, and
```

R51 is hydrogen;

or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

13. (Currently amended) Compounds A compound according to claim 1 any of the claims 1 to 10, wherein said compounds have the formula I,

in which

either

R1 is halogen, nitro, amino, mono- or di-1-4C-alkylamino, 1-4C-alkyl, 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, 3-7C-cycloalkoxy, 3-7C-cycloalkylmethoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy, and

R2 is 1-4C-alkoxy,

or

R1 is 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, 3-7C-cycloalkoxy, 3-7C-cycloalkylmethoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy, and

R2 is halogen,

and

R3 is 1-4C-alkoxy,

and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1 a]isoquinoline
pyrrolo[2,1-a]isoquinoline
ring;

and in which

either

R4 is hydrogen,

R41 is hydrogen,

R5 is hydrogen, and

R51 is hydrogen,

or

R4 is 1-4C-alkyl,

R41 is hydrogen or 1-4C-alkyl,

R5 is hydrogen, and

R51 is hydrogen;

or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

14. (Currently amended) Compounds A compound according to claim 1 any of the claims 1 to 10, wherein said compounds have the formula I, in which

- R1 is halogen, 1-4C-alkoxy, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy, such as e.g. chlorine, methoxy, 2 methoxy ethoxy or difluoromethoxy,
- R2 is 1-4C-alkoxy, such-as-e-g. methoxy,
- R3 is 1-4C-alkoxy, such as e.g. methoxy,
- and none of R1, R2 and R3 is bound to the 10-position of the pyrrolo[2.1-a]isoquinoline
  pyrrolo[2,1-a]isoquinoline ring,

and

R4 is hydrogen, or 1-4C-alkyl such as e.g. methyl,

R41 is hydrogen, or 1-4C-alkyl such as e.g. methyl,

R5 is hydrogen, and

R51 is hydrogen;

or

- R1 is halogen, 1-4C-alkoxy-2-4C-alkoxy, or completely or predominantly fluorine-substituted 1-4C-alkoxy, such as e.g. chlorine, 2 methoxy ethoxy or difluoromethoxy,
- R2 is 1-4C-alkoxy, such as e.g. methoxy,
- R3 is hydrogen,

```
and none of R1, R2 and R3 is bound to the 10-position of
    the
            pyrrolo[2.1 a] isoquinoline
                                            pyrrolo[2,1-a]-
    isoquinoline ring,
and
R4
    is hydrogen, or 1-4C-alkyl such as e.g. methyl,
R41 is hydrogen, or 1-4C-alkyl such as e.g. methyl,
R5 is hydrogen, and
R51 is hydrogen;
or
R1
   is 1-4C-alkoxy, such as e.g. methoxy,
R2
    is 1-4C-alkoxy, such as e.g. methoxy,
   is hydrogen,
R3
and none of R1, R2 and R3 is bound to the 10-position of
            pyrrolo[2.1 a] isoquinoline
pyrrolo[2,1-a] -
   the
   isoquinoline ring,
and
R4 is 1-4C-alkyl, such as-e.g. methyl,
R41 is hydrogen, or 1-4C-alkyl such as e.g. methyl,
R5 is hydrogen, and
R51 is hydrogen;
and
R6 is 1-4C-alkyl, such as e.g. methyl;
```

- or a salt, stereoisomer, hydrate or hydrate of a salt thereof
- and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.
- 15. (Currently amended) Compounds A compound according to  $\frac{\text{claim 1}}{\text{claim 5}} \frac{\text{any of the claims 1 to 10}}{\text{to 10}}$ , wherein said compounds have the formula I,

in which

- R1 is fluorine, chlorine, 1-2C-alkoxy-2-3C-alkoxy, or completely or predominantly fluorine-substituted 1-2C-alkoxy, such as e.g. chlorine, 2-methoxy-or difluoromethoxy,
- R2 is 1-2C-alkoxy, such as e.g. methoxy,
- R3 is hydrogen,
- and none of R1 and R2 is bound to the 7- or 10-position of the pyrrolo[2.1 a] isoquinoline ring,

and

R4 is hydrogen,

R41 is hydrogen,

R5 is hydrogen, and

R51 is hydrogen;

or

R1 is 1-2C-alkoxy, such as e.g. methoxy,

R2 is 1-2C-alkoxy, such as e.g. methoxy,

R3 is hydrogen,

and none of R1 and R2 is bound to the 7- or 10-position of the pyrrolo[2.1-a]isoquinoline
pyrrolo[2,1-a]isoquinoline ring,

and

R4 is 1-2C-alkyl, such as e.g. methyl,

R41 is hydrogen, or 1-2C-alkyl such as e.g. methyl,

R5 is hydrogen, and

R51 is hydrogen;

or

R1 is fluorine, chlorine, 1-2C-alkoxy-2-3C-alkoxy, or completely or predominantly fluorine-substituted 1-2C-alkoxy, such as e.g. chlorine, 2-methoxy-ethoxy or difluoromethoxy,

R2 is 1-2C-alkoxy, such as e.g. methoxy,

R3 is hydrogen,

and none of R1 and R2 is bound to the 7- or 10-position of the pyrrolo[2.1-a]isoquinoline
pyrrolo[2,1-a]isoquinoline ring,

and

R4 is 1-2C-alkyl, such as e.g. methyl,

R41 is hydrogen, or 1-2C-alkyl such-as e.g. methyl,

R5 is hydrogen, and

R51 is hydrogen;

and

R6 is 1-2C-alkyl, such as e.g. methyl;

or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

16. (Currently amended) Compounds A compound according to  $\frac{\text{claim 1}}{\text{claim 5}} \frac{\text{any of the claims 1 to 10}}{\text{to 10}}$ , wherein said compounds have the formula I,

in which

R1 is fluorine, chlorine, 1-2C-alkoxy-2-3C-alkoxy, or completely or predominantly fluorine-substituted 1-2C-alkoxy, such as e.g. chlorine, 2 methoxy ethoxy or difluoromethoxy,

- R2 is 1-2C-alkoxy, such as e.g. methoxy,
- R3 is hydrogen,
- and none of R1 and R2 is bound to the 7- or 10-position of the pyrrolo[2.1 a] isoquinoline ring,

and

R4 is 1-2C-alkyl, such as e.g. methyl,

R41 is hydrogen, or 1-2C-alkyl such as e.g. methyl,

R5 is hydrogen, and

R51 is hydrogen;

or a salt, stereoisomer, hydrate or hydrate of a salt thereof

and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.

17. (Currently amended) Compounds A compound according to claim 1 any of the preceding claims, wherein said compounds have the formula I,

in which

R6 is 1-4C-alkyl, such as e.q. methyl,

R7 is Het2, 4-hydroxy-3,5-dimethylphenyl, 3-dimethylaminophenyl, or naphthyl, in which

- Het2 is a monocyclic or fused bicyclic 5- to 10-membered heteroaryl radical comprising one to three heteroatoms, each of which is selected from [[a]] the group consisting of nitrogen, oxygen and sulfur, such as, for example, quinolyl, e.g. quinolin-4-yl;
- R8 is -C(0) -OR9, in which
- R9 is 1-4C-alkyl, such as e.g. ethyl,
- or a salt, stereoisomer, hydrate or hydrate of a salt thereof
- and the salts, stereoisomers, hydrates and hydrates of the salts of these compounds.
- 18. (Currently amended) A compound according to claim 1, wherein said compound is selected from the group consisting of:
- 1. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3,6,6trimethyl-5,6-dihydro-pyrrolo[2,1-α]isoquinoline-1carboxylic acid ethyl ester,
- 2. (6RS)-2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3,6-dimethyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,

- 3. (6RS)-8,9-Dimethoxy-3,6-dimethyl-2-(3,4,5-trimethoxy-phenyl)-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 4. 9-(1,1-Difluoro-methoxy)-2-(3-dimethylamino-phenyl)-8methoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 9-(1,1-Difluoro-methoxy)-2-(4-hydroxy-3,5-dimethyl-phenyl)-8-methoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]-isoquinoline-1-carboxylic acid ethyl ester,
- 6. 9-(1,1-Difluoro-methoxy)-8-methoxy-3-methyl-2-(3,4,5-trimethoxy-phenyl)- 5,6-dihydro-pyrrolo[2,1-a]-isoquinoline-1-carboxylic acid ethyl ester,
- 7. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9,10-trimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 8. 8-(1,1-Difluoro-methoxy)-9-methoxy-3-methyl-2-(3,4,5-trimethoxy-phenyl)-5,6-dihydro-pyrrolo[2,1-a]-isoquinoline-1-carboxylic acid ethyl ester,
- 9. 8-(1,1-Difluoro-methoxy)-2-(4-hydroxy-3,5-dimethyl-phenyl)-9-methoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]-isoquinoline-1-carboxylic acid ethyl ester,

- 10. 8-(1,1-Difluoro-methoxy)-2-(3-dimethylamino-phenyl)-9methoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 11. 8,9-(1,1-Difluoro-methylenedioxy)-2-(3-dimethylamino-phenyl)-3-methyl-5,6-dihydro-pyrrolo[2,1-a]-isoquinoline-1-carboxylic acid ethyl ester,
- 12. 8,9-(1,1-Difluoro-methylenedioxy)-2-(4-hydroxy-3,5-dimethyl-phenyl)-3-methyl-5,6-dihydro-pyrrolo[2,1-a]-isoquinoline-1-carboxylic acid ethyl ester,
- 13. 8,9-(1,1-Difluoro-methylenedioxy)-3-methyl-2-(3,4,5-trimethoxy-phenyl)-5,6-dihydro-pyrrolo[2,1-a]-isoquinoline-1-carboxylic acid ethyl ester,
- 14. 9-Chloro-2-(4-hydroxy-3,5-dimethyl-phenyl)-8-methoxy-3methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1carboxylic acid ethyl ester,
- 9-Chloro-8-methoxy-3-methyl-2-(3,4,5-trimethoxy-phenyl)-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 9-Chloro-2-(3-dimethylamino-phenyl)-8-methoxy-3-methyl5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic
  acid ethyl ester,

- 17. 8-Chloro-2-(4-hydroxy-3,5-dimethyl-phenyl)-9-methoxy-3methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1carboxylic acid ethyl ester,
- 18. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-9-methoxy-8-(2-methoxy-ethoxy)-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 19. 9-Methoxy-8-(2-methoxy-ethoxy)-3-methyl-2-(3,4,5-trimethoxy-phenyl)-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 9-Methoxy-8-(2-methoxy-ethoxy)-3-methyl-2-naphthalen-1yl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic
  acid ethyl ester,
- 9-Fluoro-2-(4-hydroxy-3,5-dimethyl-phenyl)-8-methoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 22. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8-methoxy-3-methyl-9-nitro-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 23. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8-methoxy-3,9-dimethyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,

- 24. 8,9-Dimethoxy-3-(2-methoxycarbonyl-ethyl)-6,6-dimethyl-2-quinolin-4-yl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethyl ester,
- 25. 9-Amino-2-(4-hydroxy-3,5-dimethyl-phenyl)-8-methoxy-3methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1carboxylic acid ethyl ester,
- 26. 1-[2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinolin-1-yl]-1-phenyl-methanone,
- 27. 4-(8,9-Dimethoxy-3-methyl-1-phenyl-5,6-dihydro-pyrrolo[2,1-a]isoquinolin-2-yl)-2,6-dimethyl-phenol,
- 28. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid cyclohexyl amide,
- 29. 1-[2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinolin-1-yl]-1-pyrrollidin-1-yl-methanone,
- 30. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid isopropylamide,
- 31. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid dimethylamide,

- 32. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid methylamide,
- 33. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid amide,
- 34. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid phenylamide,
- 35. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid ethylamide,
- 36. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid sec-butylamide, and
- 37. 2-(4-Hydroxy-3,5-dimethyl-phenyl)-8,9-dimethoxy-3-methyl-5,6-dihydro-pyrrolo[2,1-a]isoquinoline-1-carboxylic acid cyclopropylamide;
- and the salts, stereoisomers, hydrates, and hydrates of the salts thereof
- or a salt, stereoisomer, hydrate or hydrate of a salt thereof.

19-20. (Canceled)

- 21. (Currently amended) A pharmaceutical composition comprising one or more compounds according to claim 1, or a pharmaceutically acceptable salt, stereoisomer, hydrate or hydrate of a salt thereof, together with a customary pharmaceutical excipient and/or vehicle excipients and/or vehicles.
- 22. (Currently amended) A method for treating hyperproliferative diseases of benign or malignant behaviour and/or disorders responsive to the induction of apoptosis, such as e.g. cancer, in a patient comprising administering to said patient a therapeutically effective amount of a compound according to claim 1, or a pharmaceutically acceptable salt, stereoisomer, hydrate or hydrate of a salt thereof.
- 23. (New) The method according to claim 22, wherein said hyperproliferative disease of benign or malignant behavior and/or disorder responsive to the induction of apoptosis is cancer.